This Page Is Inserted by IFW Operations and is not a part of the Official Record

BEST AVAILABLE IMAGES

Defective images within this document are accurate representations of the original documents submitted by the applicant.

Defects in the images may include (but are not limited to):

- BLACK BORDERS
- TEXT CUT OFF AT TOP, BOTTOM OR SIDES
- FADED TEXT
- ILLEGIBLE TEXT
- SKEWED/SLANTED IMAGES
- COLORED PHOTOS
- BLACK OR VERY BLACK AND WHITE DARK PHOTOS
- GRAY SCALE DOCUMENTS

IMAGES ARE BEST AVAILABLE COPY.

As rescanning documents will not correct images, please do not report the images to the Image Problem Mailbox.



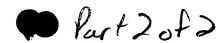
WEST

Freeform Search

Database:	US Patents Full-Text Database US Pre-Grant Publication Full-Text Database JPO Abstracts Database EPO Abstracts Database Derwent World Patents Index IBM Technical Disclosure Bulletins					
Term:						
Display: Generate:	Documents in <u>Display Format</u> : CIT, K Starting with Number 1 ○ Hit List ● Hit Count ○ Image					
	Search Clear Help Logout Interrupt					
	Main Menu Show S Numbers Edit S Numbers Preferences					
Search History						

Today's Date: 9/24/2001

DB Name	Query	Hit Count	Set Name
USPT	tangeretin	17	<u>L4</u>
USPT	tangeritin	1	<u>L3</u>
USPT	5792448	2	<u>L2</u>
EPAB,DWPI	2687572	5	L1



Trying 3106016892...Open

Welcome to STN International! Enter x:x
LOGINID:ssspta1621mxw
PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * * SESSION RESUMED IN FILE 'STNGUIDE' AT 11:04:25 ON 24 SEP 2001 FILE 'STNGUIDE' ENTERED AT 11:04:25 ON 24 SEP 2001 COPYRIGHT (C) 2001 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.00 158.39 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.76

=> d his

(FILE 'HOME' ENTERED AT 10:46:59 ON 24 SEP 2001)

FILE 'REGISTRY' ENTERED AT 10:47:09 ON 24 SEP 2001

L1 STRUCTURE UPLOADED

FILE 'REGISTRY' ENTERED AT 10:50:11 ON 24 SEP 2001 L4 1 S L3

FILE 'CAPLUS' ENTERED AT 10:51:31 ON 24 SEP 2001

L5 369 S L3

L6 3 S L5 AND COSMETIC

L7 53 S L3/THU

L8 5 S L7 AND SKIN

L9 17 S L3/PUR

FILE 'STNGUIDE' ENTERED AT 10:57:17 ON 24 SEP 2001

=> file registry COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.00 158.39 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.76

FILE 'REGISTRY' ENTERED AT 11:04:38 ON 24 SEP 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2001 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8 DICTIONARY FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8

TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

Please note that search-term pricing does apply when

conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=> d L3/rn

'L3' MUST END IN '/Q', '/A', '/L', '/S' OR '/B'

The saved name for a query (or structure or screen set) must end with $^\prime/Q^\prime$. The saved name for an answer set must end with $^\prime/A^\prime$. The saved name for an L# list must end with $^\prime/L^\prime$. SDI request names must end with $^\prime/S^\prime$. To see a list of all saved query, answer set,, and L# list names for this loginid, enter "DISPLAY SAVED" at an arrow prompt (=>). Enter "DISPLAY SAVED/S" to see a list of SDI request names. Enter "DISPLAY SAVED/B" to see a list of BATCH search requests.

=> d L3

L3 ANSWER 1 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 315674-37-4 REGISTRY

CN 4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)-, labeled with tritium (9CI) (CA INDEX NAME)

MF C20 H20 O7

SR CAS Registry Services

IL XH-3

=> d L3 2-14

L3 ANSWER 2 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 229152-82-3 REGISTRY

CN .beta.-Cyclodextrin, compd. with

5, 6, 7, 8-tetramethoxy-2-(4-methoxyphenyl)-

4H-1-benzopyran-4-one (1:1) (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)-, compd. with .beta.-cyclodextrin (1:1) (9CI)

FS STEREOSEARCH

MF C42 H70 O35 . C20 H20 O7

SR CA

LC STN Files: CA, CAPLUS

CM 1

CRN 7585-39-9 CMF C42 H70 O35

Absolute stereochemistry.

PAGE 1-A

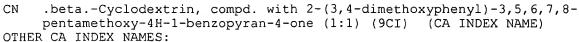
PAGE 2-A

CM2

CRN 481-53-8 CMF C20 H20 O7

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L3 ANSWER 3 OF 14 REGISTRY COPYRIGHT 2001 ACS RN 229152-81-2 REGISTRY



CN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5,6,7,8-pentamethoxy-, compd. with .beta.-cyclodextrin (1:1) (9CI)

FS STEREOSEARCH

MF C42 H70 O35 . C22 H24 O9

SR CA

LC STN Files: CA, CAPLUS

CM 1

CRN 7585-39-9 CMF C42 H70 O35

Absolute stereochemistry.

PAGE 1-A

ÖН

PAGE 2-A

CRN 1178-24-1 CMF C22 H24 O9

2

CM

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L3 ANSWER 4 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 229152-80-1 REGISTRY

CN .beta.-Cyclodextrin, compd. with 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-4H-1-benzopyran-4-one (1:1) (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES:

CN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-, compd. with .beta.-cyclodextrin (1:1) (9CI)

FS STEREOSEARCH

MF C42 H70 O35 . C21 H22 O8

SR CA

LC STN Files: CA, CAPLUS

CM 1

CRN 7585-39-9 CMF C42 H70 O35

Absolute stereochemistry.

PAGE 2-A

CM 2

CRN 478-01-3 CMF C21 H22 O8

1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L3 ANSWER 5 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 52412-94-9 REGISTRY

CN 4H-1-Benzopyran-4-one, 3,5,6,7,8-pentamethoxy-2-[4-(methoxy-d3)phenyl]-(9CI) (CA INDEX NAME)

MF C21 H19 D3 O8

LC STN Files: CA, CAPLUS

1 REFERENCES IN FILE CA (1967 TO DATE) 1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L3 ANSWER 6 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 51901-05-4 REGISTRY

CN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-, mixt. with sulfinylbis[methane] (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Methane, sulfinylbis-, mixt. contg. (9CI)

OTHER NAMES:

CN Dimethyl sulfoxide-Nobiletine mixture

MF C21 H22 O8 . C2 H6 O S

CI MXS

LC STN Files: CA, CAPLUS, TOXLIT

CM 1

CRN 478-01-3 CMF C21 H22 O8

CM 2

CRN 67-68-5 CMF C2 H6 O S

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L3 ANSWER 7 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 34170-18-8 REGISTRY

CN 4H-1-Benzopyran-4-one, 3,5,6,7,8-pentamethoxy-2-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Flavone, 3,4',5,6,7,8-hexamethoxy- (6CI, 8CI)

MF C21 H22 O8

LC STN Files: BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, TOXLIT (*File contains numerically searchable property data)

- 9 REFERENCES IN FILE CA (1967 TO DATE)
- 9 REFERENCES IN FILE CAPLUS (1967 TO DATE)
- 1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L3 ANSWER 8 OF 14 REGISTRY COPYRIGHT 2001 ACS

```
24027-55-2 REGISTRY
RN
     4H-1-Benzopyran-4-one, 3,5,7,8-tetramethoxy-2-(4-methoxyphenyl)- (9CI)
CN
     (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Flavone, 3,4',5,7,8-pentamethoxy- (7CI, 8CI)
OTHER NAMES:
     3,5,7,8,4'-Pentamethoxyflavone
CN
     Tambulin dimethyl ether
CN
FS
     3D CONCORD
MF
     C20 H20 O7
                  BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, TOXLIT
LC
     STN Files:
         (*File contains numerically searchable property data)
       OMe
```

50 REFERENCES IN FILE CA (1967 TO DATE) 50 REFERENCES IN FILE CAPLUS (1967 TO DATE)

```
ANSWER 10 OF 14 REGISTRY COPYRIGHT 2001 ACS
L3
RN
     7741-47-1 REGISTRY
CN
     4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5,7,8-tetramethoxy-
(9CI)
      (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Flavone, 3,3',4',5,7,8-hexamethoxy- (6CI, 7CI, 8CI)
OTHER NAMES:
CN
     3,3',4',5,7,8-Hexamethoxyflavone
     3,5,7,8,3',4'-Hexamethoxyflavone
CN
     Gossypetin hexamethyl ether
CN
FS
     3D CONCORD
MF
     C21 H22 O8
LC
     STN Files:
                  AGRICOLA, BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, NAPRALERT,
       SPECINFO, TOXLINE, TOXLIT
         (*File contains numerically searchable property data)
       OMe
MeO
                         OMe
               OMe
       OMe
                    OMe
              33 REFERENCES IN FILE CA (1967 TO DATE)
              33 REFERENCES IN FILE CAPLUS (1967 TO DATE)
               6 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
L3
     ANSWER 11 OF 14 REGISTRY COPYRIGHT 2001 ACS
RN
     6601-66-7 REGISTRY
CN
     4H-1-Benzopyran-4-one, 5,7,8-trimethoxy-2-(4-methoxyphenyl)- (9CI)
     INDEX NAME)
OTHER CA INDEX NAMES:
     Flavone, 4',5,7,8-tetramethoxy- (7CI, 8CI)
CN
OTHER NAMES:
CN
     4',5,7,8-Tetramethoxyflavone
CN
     6-Demethoxytangeretin
CN
     6-Demethoxytangeritin
CN
     Tetra-O-methylisoscutellarein
FS
     3D CONCORD
     C19 H18 O6
MF
LC
     STN Files:
                  AGRICOLA, BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD,
CAPLUS,
       NAPRALERT, SPECINFO, TOXLIT
         (*File contains numerically searchable property data)
       OMe
MeO
                        OMe
```

OMe

0

```
1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
               46 REFERENCES IN FILE CAPLUS (1967 TO DATE)
               1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
     ANSWER 12 OF 14 REGISTRY COPYRIGHT 2001 ACS
     1178-24-1 REGISTRY
RN
     4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5,6,7,8-pentamethoxy-
CN
             (CA INDEX NAME)
     (9CI)
OTHER CA INDEX NAMES:
     Flavone, 3,3',4',5,6,7,8-heptamethoxy- (6CI, 7CI, 8CI)
     3', 4', 3, 5, 6, 7, 8-Heptamethoxyflavone
CN
     3, 3', 4', 5, 6, 7, 8-Heptamethoxyflavone
     3, 5, 6, 7, 8, 3', 4'-Heptamethoxyflavone
CN
MF
     C22 H24 O9
CI
     COM
                  AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, CA,
LC
     STN Files:
       CANCERLIT, CAOLD, CAPLUS, CHEMCATS, IFICDB, IFIPAT, IFIUDB, MEDLINE,
       NAPRALERT, TOXLINE, TOXLIT, USPATFULL
         (*File contains numerically searchable property data)
       OMe
MeO
MeO
                         0Me
              OMe
       OMe
            0
                    OMe
             103 REFERENCES IN FILE CA (1967 TO DATE)
             103 REFERENCES IN FILE CAPLUS (1967 TO DATE)
               9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
L3
     ANSWER 13 OF 14 REGISTRY COPYRIGHT 2001 ACS
RN
     481-53-8 REGISTRY
     4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)- (9CI)
CN
     (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Flavone, 4',5,6,7,8-pentamethoxy- (7CI, 8CI)
     Tangeretin (6CI)
OTHER NAMES:
CN
     4',5,6,7,8-Pentamethoxyflavone
CN
     5, 6, 7, 8, 4'-Pentamethoxyflavone
CN
     Ponkanetin
CN
     Tangeritin
MF
     C20 H20 O7
CI
     COM
LC
                  ADISINSIGHT, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CHEMCATS, CHEMLIST,
       CSCHEM, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE,
       NAPRALERT, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
                      EINECS**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

46 REFERENCES IN FILE CA (1967 TO DATE)

271 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

271 REFERENCES IN FILE CAPLUS (1967 TO DATE)

14 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L3 ANSWER 14 OF 14 REGISTRY COPYRIGHT 2001 ACS

RN 478-01-3 REGISTRY

CN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-(9CI)

(CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Flavone, 3',4',5,6,7,8-hexamethoxy- (7CI, 8CI)

CN Nobiletin (6CI)

OTHER NAMES:

CN 3', 4', 5, 6, 7, 8-Hexamethoxyflavone

CN 5, 6, 7, 8, 3', 4'-Hexamethoxyflavone

MF C21 H22 O8

CI COM

LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMCATS, CSCHEM,
DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, NAPRALERT,
PROMT, SPECINFO, TOXLINE, TOXLIT, USPATFULL
(*File contains numerically searchable property data)

210 REFERENCES IN FILE CA (1967 TO DATE)

210 REFERENCES IN FILE CAPLUS (1967 TO DATE)

14 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

Part lof 2

Trying 3106016892...Open

Welcome to STN International! Enter x:x

LOGINID:ssspta1621mxw

PASSWORD:

t == *

TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS	1			Web Page URLs for STN Seminar Schedule - N. America
NEWS	2	Dec	17	The CA Lexicon available in the CAPLUS and CA files
NEWS	3	Feb	06	Engineering Information Encompass files have new names
NEWS	4	Feb	16	TOXLINE no longer being updated
NEWS	5	Apr	23	Search Derwent WPINDEX by chemical structure
NEWS	6	Apr	23	PRE-1967 REFERENCES NOW SEARCHABLE IN CAPLUS AND CA
NEWS	7	May	07	DGENE Reload
NEWS	8	Jun	20	Published patent applications (A1) are now in USPATFULL
NEWS	9	JUL	13	New SDI alert frequency now available in Derwent's DWPI and DPCI
NEWS	10	Aug	23	<pre>In-process records and more frequent updates now in MEDLINE</pre>
NEWS	11	Aug	23	PAGE IMAGES FOR 1947-1966 RECORDS IN CAPLUS AND CA
NEWS	12	Aug	23	Adis Newsletters (ADISNEWS) now available on STN
NEWS	13	Sep	17	IMSworld Pharmaceutical Company Directory name change to PHARMASEARCH
				CO I HANNADEANON
NEWS EXPRESS			gust 15 CURRENT WINDOWS VERSION IS V6.0c,	
				RRENT MACINTOSH VERSION IS V6.0 (ENG) AND V6.0J (JP),
NEELO				D CURRENT DISCOVER FILE IS DATED 07 AUGUST 2001
NEWS		-		N Operating Hours Plus Help Desk Availability
NEWS	INTE	ΣR	Gei	neral Internet Information
NEWS	WS LOGIN Welcome Banner and News Items			

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

NEWS PHONE Direct Dial and Telecommunication Network Access to STN

CAS World Wide Web Site (general information)

FILE 'HOME' ENTERED AT 10:46:59 ON 24 SEP 2001

=> file registry COST IN U.S. DOLLARS

NEWS WWW

SINCE FILE TOTAL ENTRY SESSION 0.15 0.15

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 10:47:09 ON 24 SEP 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2001 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8 DICTIONARY FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8

TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>

Uploading 09837129a.str

L1 STRUCTURE UPLOADED

STR

=> d L1

L1 HAS NO ANSWERS

L1

Structure attributes must be viewed using STN Express query preparation.

=> s L1 sss sam

SAMPLE SEARCH INITIATED 10:47:50 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 255 TO ITERATE

100.0% PROCESSED 255 ITERATIONS SEARCH TIME: 00.00.01

1 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COM

CTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 4142 TO 6058 PROJECTED ANSWERS: 1 TO 80

```
L2 1 SEA SSS SAM L1
```

=> d L1 scan

L1 HAS NO ANSWERS

=> d 12 scan

L2 1 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)- (9CI)

MF C20 H20 O7

CI COM

ALL ANSWERS HAVE BEEN SCANNED

```
=> d L1 sss full
L1 HAS NO ANSWERS
'SSS FULL ' IS NOT A VALID STRUCTURE FORMAT KEYWORD
Structure Formats
SIA ---- Structure Image, Attributes, and map table if it contains
          data. (Default)
SIM ---- Structure IMage.
SAT ---- Structure ATtributes and map table if it contains data.
SCT ---- Structure Connection Table and map table if it contains
SDA ---- All Structure DAta (image, attributes, connection table and
          map table if it contains data).
NOS ---- NO Structure data.
ENTER STRUCTURE FORMAT (SIM), NOS:
ENTER STRUCTURE FORMAT (SIM), NOS:0
'0' IS NOT A VALID STRUCTURE FORMAT KEYWORD
Structure Formats
SIA ---- Structure Image, Attributes, and map table if it contains
          data. (Default)
SIM ---- Structure IMage.
SAT ---- Structure ATtributes and map table if it contains data.
SCT ---- Structure Connection Table and map table if it contains
SDA ---- All Structure DAta (image, attributes, connection table and
         map table if it contains data).
NOS ---- NO Structure data.
ENTER STRUCTURE FORMAT (SIM), NOS:sim
L1
                STR
```

Structure attributes must be viewed using STN Express query preparation.

=> s L1 sss full FULL SEARCH INITIATED 10:49:04 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 5289 TO ITERATE

100.0% PROCESSED 5289 ITERATIONS

SEARCH TIME: 00.00.01

L3 14 SEA SSS FUL L1

=> d L3 scan

G1 MeO, H

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 3,5,6,7,8-pentamethoxy-2-(4-methoxyphenyl)- (9CI) MF C21 H22 O8

14 ANSWERS

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN .beta.-Cyclodextrin, compd. with 2-(3,4-dimethoxyphenyl)-5,6,7,8tetramethoxy-4H-1-benzopyran-4-one (1:1) (9CI)
MF C42 H70 O35 . C21 H22 O8

CM 1

Absolute stereochemistry.

PAGE 1-A

CM 2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)-, labeled with tritium (9CI)

MF C20 H20 O7

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5,6,7,8-pentamethoxy(9CI)

MF C22 H24 O9

CI COM

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,7,8-trimethoxy- (9CI)

MF C20 H20 O7

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-, mixt. with sulfinylbis[methane] (9CI)

MF C21 H22 O8 . C2 H6 O S

CI MXS

> CM 1

CM2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN .beta.-Cyclodextrin, compd. with 2-(3,4-dimethoxyphenyl)-3,5,6,7,8pentamethoxy-4H-1-benzopyran-4-one (1:1) (9CI) C42 H70 O35 . C22 H24 O9

MF

CM1

Absolute stereochemistry.

PAGE 1-A

PAGE 2-A

CM 2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphen

IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-5,6,7,8-tetramethoxy-(9CI)

MF C21 H22 O8

CI COM

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 5,7,8-trimethoxy-2-(4-methoxyphenyl)- (9CI)

MF C19 H18 O6

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 3,5,7,8-tetramethoxy-2-(4-methoxyphenyl)- (9CI)

MF C20 H20 O7

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

MF C21 H19 D3 O8

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS IN .beta.-Cyclodextrin, compd. with 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)-

4H-1-benzopyran-4-one (1:1) (9CI)

MF C42 H70 O35 . C20 H20 O7

CM 1

Absolute stereochemistry.

PAGE 1-A

PAGE 2-A

CM 2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 4H-1-Benzopyran-4-one, 5,6,7,8-tetramethoxy-2-(4-methoxyphenyl)- (9CI)

MF C20 H20 O7

CI COM

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN 4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5,7,8-tetramethoxy-(9CI)

MF C21 H22 O8

ALL ANSWERS HAVE BEEN SCANNED

=> file registry
COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL
ENTRY SESSION
134.80 134.95

FILE 'REGISTRY' ENTERED AT 10:50:11 ON 24 SEP 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2001 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8 DICTIONARY FILE UPDATES: 23 SEP 2001 HIGHEST RN 358331-51-8

TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=> d his

(FILE 'HOME' ENTERED AT 10:46:59 ON 24 SEP 2001)

FILE 'REGISTRY' ENTERED AT 10:47:09 ON 24 SEP 2001

L1 STRUCTURE UPLOADED

FILE 'REGISTRY' ENTERED AT 10:50:11 ON 24 SEP 2001

=> s L3

SAMPLE SEARCH INITIATED 10:51:13 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 255 TO ITERATE

100.0% PROCESSED 255 ITERATIONS 1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

PROJECTED ITERATIONS: 4142 TO 6058

PROJECTED ANSWERS: 1 TO 80

L4 1 SEA SSS SAM L1

=> file caplus

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
0.62 135.57

FILE 'CAPLUS' ENTERED AT 10:51:31 ON 24 SEP 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2001 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications.

FILE COVERS 1947 - 24 Sep 2001 VOL 135 ISS 14 FILE LAST UPDATED: 23 Sep 2001 (20010923/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

CAplus now provides online access to patents and literature covered in CA from 1947 to the present. On April 22, 2001, bibliographic information and abstracts were added for over 2.2 million references published in CA from 1947 to 1966.

The CA Lexicon is now available in the Controlled Term (/CT) field. Enter HELP LEXICON for full details.

Attention, the CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

=> s L3

L5 369 L3

=> s L5 and cosmetic

37070 COSMETIC 35434 COSMETICS 48574 COSMETIC

(COSMETIC OR COSMETICS)

L6 3 L5 AND COSMETIC

=> d L6 1-3 ti

L6 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2001 ACS

TI Lipolytic agent as health food and **cosmetic** for treatment of obesity

L6 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2001 ACS

TI **Cosmetic** composition containing a metal oxide nanopigment and an antioxidant for hair and skin protection

L6 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2001 ACS

TI Cosmetic compositions containing flavonoids

=> d L6 3 ibib, abs

L6 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2001 ACS ACCESSION NUMBER: 1994:173131 CAPLUS

DOCUMENT NUMBER: 120:173131

TITLE: Cosmetic compositions containing flavonoids INVENTOR(S): Beck, Irena; Hocquaux, Michel; Tournaire, Cecile

PATENT ASSIGNEE(S): Oreal S. A., Fr. SOURCE: Fr. Demande, 14 pp.

CODEN: FRXXBL

DOCUMENT TYPE: Patent LANGUAGE: French

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE FR 1992-2247 19920226 FR 2687572 A1 19930827 B1 19950609 FR 2687572 OTHER SOURCE(S): MARPAT 120:173131 Cosmetic compns. contg. flavonoids which have a double bond on position 2,3 and OH on position 3 are used for protection of the skin (Markush structure given). A cream contained Sinowax AO 7, Gelcol 2, cetyl alc. 1.5, iso-Pr myristate 1.5, vaseline 15.0, propylene glycol 10.0, glycerin 10.0, catechin 0.5, preservatives q.s. and water q.s. 100g. => s L3/thu 369 L3 396285 THU/RL L7 53 L3/THU (L3 (L) THU/RL)=> s L7 and skin 156526 SKIN 5617 SKINS 159382 SKIN (SKIN OR SKINS) 5 L7 AND SKIN L8 => d L8 1-5 ti 1.8 ANSWER 1 OF 5 CAPLUS COPYRIGHT 2001 ACS TICompositions and methods for treatment of neoplastic diseases with combinations of limonoids, flavonoids and tocotrienols L8ANSWER 2 OF 5 CAPLUS COPYRIGHT 2001 ACS TΙ Cancer chemopreventive activity of 3,5,6,7,8,3',4'-heptamethoxyflavone from the peel of citrus plants ANSWER 3 OF 5 CAPLUS COPYRIGHT 2001 ACS L8 TΙ Inhibitory effect of citrus nobiletin on phorbol ester-induced skin inflammation, oxidative stress, and tumor promotion in mice L8 ANSWER 4 OF 5 CAPLUS COPYRIGHT 2001 ACS TΙ Inhibitory effect of flavonoids from Citrus plants on Epstein-Barr virus activation and two-stage carcinogenesis of skin tumors 18 ANSWER 5 OF 5 CAPLUS COPYRIGHT 2001 ACS TΙ Compositions and methods of inhibiting neoplastic and cardiovascular diseases with compounds related to limocitrin and 5-desmethyl sinensetin => s L3/pur 369 L3 151393 PUR/RL L9 17 L3/PUR (L3 (L) PUR/RL) => d L9 1-17 ti L9 ANSWER 1 OF 17 CAPLUS COPYRIGHT 2001 ACS Inhibitors of 15-Lipoxygenase from Orange Peel

ANSWER 2 OF 17 CAPLUS COPYRIGHT 2001 ACS TIStudies on chemical constituents of processed green tangerine peel ANSWER 3 OF 17 CAPLUS COPYRIGHT 2001 ACS 1.9 ΤI Pharmaceuticals and foods containing flavonoids as inhibitors of formation of matrix metalloproteinase (MMP) and its precursor L9 ANSWER 4 OF 17 CAPLUS COPYRIGHT 2001 ACS A new oxyprenyl coumarin and highly methylated flavones from the exudate TΤ of Ozothamnus lycopodioides (Asteraceae) T.9 ANSWER 5 OF 17 CAPLUS COPYRIGHT 2001 ACS ΤI Chemical constituents of Tripterygium wilfordii 1.9 ANSWER 6 OF 17 CAPLUS COPYRIGHT 2001 ACS Isolation and Identification of New Polymethoxyflavonoids from Dancy TITangerine Leaves L9 ANSWER 7 OF 17 CAPLUS COPYRIGHT 2001 ACS TΙ Studies on the chemical constituents Yengmaoercao (Hedyotis lindleyanu) L9 ANSWER 8 OF 17 CAPLUS COPYRIGHT 2001 ACS Study on the constituents of flavonoids and other substances from peel of ΤI Citrus Sinensis (L.) Osbeck L9 ANSWER 9 OF 17 CAPLUS COPYRIGHT 2001 ACS ΤI The genuineness of citrus oils. Note LV. Composition of mandarin essential oil produced in 1996/97 L9 ANSWER 10 OF 17 CAPLUS COPYRIGHT 2001 ACS On the chemical, botanical, and chemotaxonomical evaluation of the genus TΙ Citrus. Part I: polymethoxyflavones of the leaf of Citrus deliciosa Ten. T. 9 ANSWER 11 OF 17 CAPLUS COPYRIGHT 2001 ACS Two New Polymethoxylated Flavones, a Class of Compounds with Potential TIAnticancer Activity, Isolated from Cold Pressed Dancy Tangerine Peel Oil Solids T.9 ANSWER 12 OF 17 CAPLUS COPYRIGHT 2001 ACS ΤI Chemical constituents from roots of Jatropha curcas 1,9 ANSWER 13 OF 17 CAPLUS COPYRIGHT 2001 ACS TΙ Isolation of anti-leukemia compounds from Citrus reticulata L9 ANSWER 14 OF 17 CAPLUS COPYRIGHT 2001 ACS Flavonoids of mandarin fruit refuse and their fungistatic action on the ΤI fungus Phoma tracheiphila L9 ANSWER 15 OF 17 CAPLUS COPYRIGHT 2001 ACS TΙ Isolation of polymethoxyflavones from sweet and Mandarin orange essential oils using column chromatography and semipreparative HPLC with a recycle valve L 9 ANSWER 16 OF 17 CAPLUS COPYRIGHT 2001 ACS ΤI Cardiotonic flavonoids from Citrus plants (Rutaceae) L9 ANSWER 17 OF 17 CAPLUS COPYRIGHT 2001 ACS

TΤ Further pyrano flavones from Neoraputia alba

=> d L9 15-16 ibib, abs

ANSWER 15 OF 17 CAPLUS COPYRIGHT 2001 ACS

ACCESSION NUMBER:

1995:538084 CAPLUS

DOCUMENT NUMBER:

122:273747

TITLE:

Isolation of polymethoxyflavones from sweet and Mandarin orange essential oils using column chromatography and semipreparative HPLC with a

recycle

valve

AUTHOR (S):

Mondello, Luigi; Dugo, Paola; Stagno d'Alcontres,

Ildefonsa

CORPORATE SOURCE:

Facolta di Farmacia, Univ. Messina, Italy Essenze, Deriv. Agrum. (1993), 63(4), 395-406

SOURCE:

CODEN: EDAGAH; ISSN: 0014-0902

DOCUMENT TYPE:

Journal

LANGUAGE:

Italian

A method for the isolation of polymethoxylated flavones (PMF) from sweet orange and mandarin essential oils using chromatog. on a silica gel column

and HPLC with recycle valve, using "peak shaving" technique is reported. The pure PMF were used to construct calibration curves against coumarin for the quant. estn. of PMF in the two essential oils.

ANSWER 16 OF 17 CAPLUS COPYRIGHT 2001 ACS

ACCESSION NUMBER:

1995:437431 CAPLUS

DOCUMENT NUMBER:

122:183243

TITLE:

Cardiotonic flavonoids from Citrus plants (Rutaceae) Itoigawa, Masataka; Takeya, Kazumi; Furukawa, Hiroshi

AUTHOR(S): CORPORATE SOURCE:

Tokaigakuen Women's College, Nagoya, 468, Japan

Biol. Pharm. Bull. (1994), 17(11), 1519-21

SOURCE:

CODEN: BPBLEO; ISSN: 0918-6158

DOCUMENT TYPE:

Journal

Ι

LANGUAGE:

English

GT

OMe OMe MeO OMe MeO OH OMe 0

AB Two flavonoids, 3,5,6,7,8,3',4'-heptamethoxyflavone (HEPTA) and natsudaidain (I) isolated from Citrus plants (Rutaceae), produced a pos. inotropic effect (PIE) on guinea-pig papillary muscle. I (pD2 4.98.+-.0.07) was more potent than HEPTA, but the max. PIE of HEPTA was greater than that of I. The PIE of HEPTA was completely inhibited by reserpinization of the guinea pig, and partially inhibited by metoprolol and carbachol. The carbachol inhibition was omitted by atropine. The mechanism of PIE of HEPTA is accounted for catecholamine release from

cardiac tissue.

=> file stng COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	22.82	158.39
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-1.76	-1.76

FILE 'STNGUIDE' ENTERED AT 10:57:17 ON 24 SEP 2001 USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT COPYRIGHT (C) 2001 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Sep 21, 2001 (20010921/UP).

=> d his

(FILE 'HOME' ENTERED AT 10:46:59 ON 24 SEP 2001)

FILE 'REGISTRY' ENTERED AT 10:47:09 ON 24 SEP 2001 STRUCTURE UPLOADED

L1 STRUCTURE UPLOAI L2 1 S L1 SSS SAM

L3 14 S L1 SSS FULL

FILE 'REGISTRY' ENTERED AT 10:50:11 ON 24 SEP 2001 L4

FILE 'CAPLUS' ENTERED AT 10:51:31 ON 24 SEP 2001

L5 369 S L3

L6 3 S L5 AND COSMETIC

L7 53 S L3/THU

L8 5 S L7 AND SKIN

L9 17 S L3/PUR

FILE 'STNGUIDE' ENTERED AT 10:57:17 ON 24 SEP 2001